

Transfer of trapped atoms between two optical tweezer potentials

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Trapped, laser-cooled rubidium atoms are transferred between two strongly focused, horizontal, orthogonally intersecting laser beams. The transfer efficiency is studied as a function of the vertical distance between the beam axes. Optimum transfer is found when the distance equals the beam waist radius. Numerical simulations reproduce well the experimental results.

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I. INTRODUCTION

The spatial control of atoms, beyond their trapping in stationary potentials, has been continuously gaining importance in investigations of ultracold gases and in the application of atomic ensembles and single atoms for cavity QED and quantum information studies. Recent progress includes the trapping and control of single atoms in dynamic potentials [1, 2], the sub-micron positioning of individual atoms with standing-wave potentials [3, 4], micro-structured and dynamic traps for Bose-Einstein condensates [5, 6] and, as another example, the realisation of chaotic dynamics in atom-optics "billiards" [7, 8].

The reaction of trapped atoms to dynamical variation of the trapping potential is one central aspect of these developments. Its understanding is important to design optimally the shape of the potential and its temporal variations which provide the desired control over the atoms and allow their manipulations. The question of efficiently steering atoms by dynamically variable light beams shares many similarities with the application of laser beams for optical tweezers, for manipulating microscopic objects such as beads or living cells [9]. In the context of quantum information processing, it is also related to recent developments towards position control of single trapped ions in complex, multi-segment ion traps [10].

We report on a particular case of manipulation of atoms by dynamic variation of light potentials formed by laser beams: the transfer of a cold cloud of rubidium atoms between two horizontal, intersecting, focused laser beams. The transfer happens by ramping the intensities in the two beams, which is done slowly enough to give all atoms time to adjust to the change. The measured signal is the amount of atoms remaining in the second beam after the first one has been fully switched off, i.e. the transfer efficiency. Our main finding is a peculiar dependence of the efficiency on the vertical distance be-

tween the two beams, which we vary within about ± 3 times the beam waist radius w_0 . The striking result is that we observe optimum transfer when the beams are not fully overlapping but when their distance is about w_0 . For the situation with maximum overlap the transfer efficiency is reduced by about a factor of two.

The observed behaviour is reproduced in numerical simulations of the situation, using classical trajectories of independent atoms. Supported by these results, our explanation is that the finite distance of w_0 between the beams is favourable for scattering an atom from one beam to the other, because it does not create any potential barrier, while at the same time the anharmonic potential mixes the degrees of freedom of orthogonal directions.

II. EXPERIMENT

In the experiment, we trap a cloud of ^{87}Rb atoms in a dipole trap formed by about 0.7 W of light from a Ti:Sapphire laser at 810.0 nm, focused to a waist diameter of about $2w_0 = 15 \mu\text{m}$. The atoms are transferred into the dipole trap from a standard magneto-optical trap (MOT) formed by three retro-reflected beams from a frequency-stabilised diode laser [11]. After switching off the MOT, the atoms are held in the dipole trap for 100 ms, during which they thermalise. Then the second, orthogonal laser beam is ramped up from 0 to 0.4 W within 100 ms, followed by ramping down the first beam to zero during the next 100 ms. After another 220 ms of waiting time, the MOT beams are switched back on, the dipole trap is switched off completely, and the fluorescence of the recaptured atoms is measured during 100 ms. The procedure is repeated several 10 times to average over shot-to-shot fluctuations in the initial number of trapped atoms.

The two dipole trapping beams are horizontal and have the same beam parameters. At their respective maximum powers of 0.7 and 0.4 W, the single-beam potentials are 150 and 86 MHz deep and have 42 and 32 kHz radial frequencies. Their ramping is done with acousto-optical modulators (AOMs), while mechanical shutters

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are used to switch them off completely before and after ramping. The beams intersect at their focal points, apart from an adjustable vertical distance which is the main experimental parameter. Adjustment happens with a precision translation stage that moves vertically the complete assembly of fiber collimator and focusing lens which provides the second trapping beam.

The main experimental result is shown in Fig. 1. On variation of the vertical distance, the transfer efficiency exhibits two maxima, around $\pm w_0$. In the symmetric situation, at maximum overlap of the beams, the efficiency is reduced, and, obviously, it goes to zero when the beam distance becomes large.

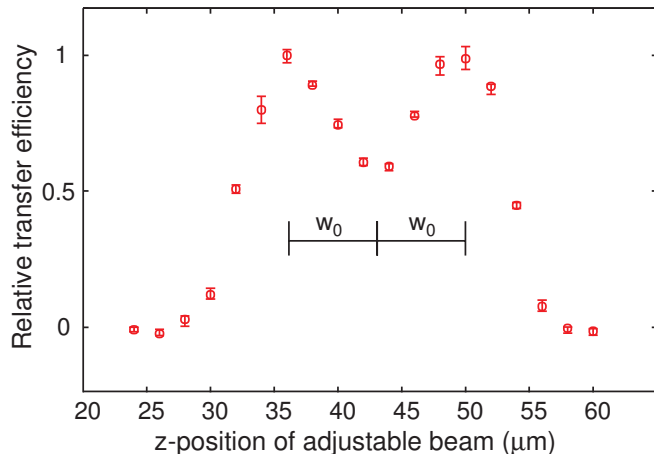


FIG. 1: Measured transfer efficiency between the two trapping beams as a function of their vertical distance. The error bars correspond to a confidence level of 68%, determined from several 10 repetitions of the same measurement. The efficiency is normalised to the maximum.

III. QUALITATIVE DESCRIPTION

The probability of an atom to be transferred from one beam to the other depends on the shape of the combined potential in the crossing region. Here we provide a qualitative discussion of the physical situation, before we present a numerical simulation of the dynamics in the next section.

We denote the three spatial directions by \hat{x} for the propagation direction of the first beam, \hat{y} for that of the second, and \hat{z} for the vertical direction. In Fig. 2, we show three relevant cases of vertical distance, assuming equal beam power. For zero distance the two beams combine to form one localised 3-dimensional potential "dimple" which is symmetric in \hat{x} and \hat{y} , and stronger in \hat{z} . For a distance larger than w_0 , a potential barrier separates the two beams. In the intermediate case of a separation by w_0 , neither a dimple nor a barrier is formed. Instead, in the \hat{z} direction the potential has a non-Gaussian, flat-bottom shape.

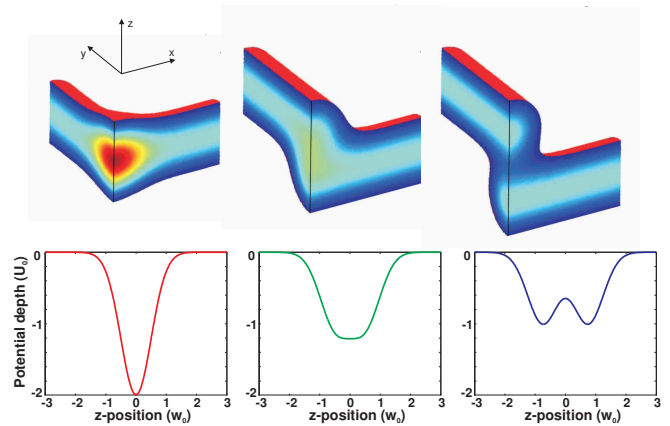


FIG. 2: Illustration of the potential in the crossing region, for the vertical distance between the beams set to 0 (left), w_0 (center) and $1.5w_0$ (right). Top: color- (or grey-scale-) coded 3-dimensional display. Bottom: vertical potential variation at the center of the crossing region (along the black line in the top display).

Some qualitative conclusions can be drawn from these pictures. At zero beam separation and for an atom with small kinetic energy, the trajectory will not be affected significantly when it traverses the beam crossing region, except for a faster radial oscillation and some longitudinal acceleration and deceleration. The potential is always close to harmonic and symmetric. Due to this symmetry, the motion will remain centered around the axis of the beam in which the atom enters the crossing region, and the three directions of motion are not mixed. Thus a transfer between the beams is not likely.

In the case of large beam separation, the potential barrier impedes the transfer, so the probability that the atom changes between the beams falls off to zero. The fall-off is expected to happen faster (i.e. at smaller beam separation) for atoms with lower energy.

The maximal transfer probability is observed in the intermediate case. We ascribe this to the asymmetric potential in the crossing region, which deflects all incoming atoms up- or downwards towards the axis of the other beam, such that the motional degrees of freedom become necessarily coupled. Moreover, the potential in the crossing region is anharmonic, which enhances the mixing between the directions.

IV. MODEL CALCULATIONS

Numerical simulations of the classical motion of a particle in the crossed-beam potential serve as a complementary approach to understand the observed dynamics. Based on the analytical expressions for the potential and its spatial gradient, the differential equations of motion were solved numerically using standard tools [12]. At typical atom temperatures of 50 μ K and densities of

several 10^{10} cm^{-3} , the classical trajectory approach is well justified. Instead of simulating the complete transfer process, which would have required significantly more computational resources, we looked at the dynamics of a single particle in the potential formed by two beams of constant, equal power. Hence these calculations highlight the effect of the coupling between the motional degrees of freedom in the crossing region, and of the potential barrier (when it exists) between the beams.

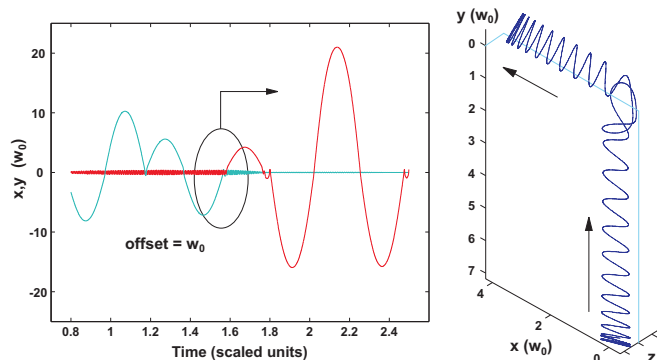


FIG. 3: Sample trajectory from the numerical simulation. Left: excursion in \hat{x} and \hat{y} , in units of w_0 , vs. time; \hat{z} -motion is not shown. Right: 3-dimensional display zooming into the circled region of the left plot which shows a transfer event.

Figure 3 shows a sample trajectory, for a displacement between the beams equal to the beam waist w_0 . The numerical calculations were programmed with high temporal resolution to account for the strongly different trap frequencies and for the anharmonic potential encountered by particles off the beam axes. In the simulations the total energy (the sum of kinetic and potential energy) was found to change slowly with time, due to accumulated numerical errors, but it was practically constant during a single transit through the crossing region, such that it served as a parameter characterising each transit event. The probability for a transfer from one beam to the other during a transit through the crossing region was then recorded as a function of the energy and the beam separation. A particle was considered to be transferred between the beams if it entered the central region from a distance larger than $3w_0$ in one beam, and left this region in the other beam to a distance larger than $3w_0$. A transit was defined as any event where the particle entered and left the crossing region in either of the beams. The transfer probability is the ratio of the number of transfers to the number of transits. The initial conditions were chosen at random; gravity was neglected [13].

Figure 4 shows a histogram of transfer probability vs. beam separation and particle energy from such a numerical investigation. The value of each bin is based on 100 to 1000 transit events.

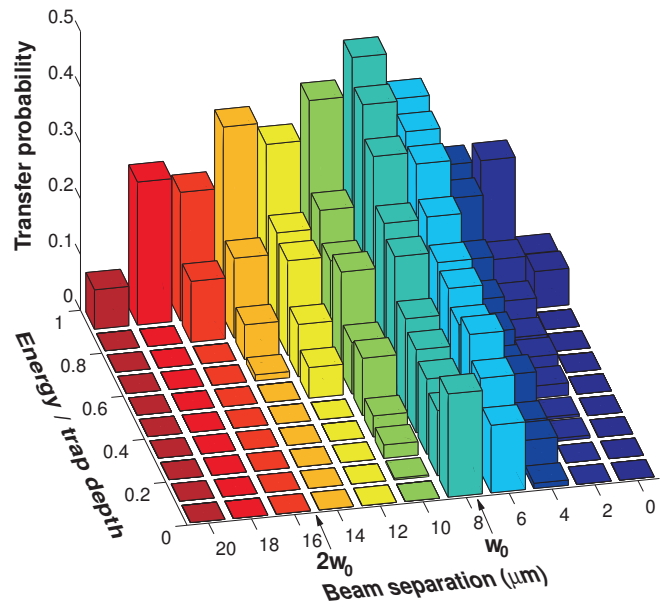


FIG. 4: Simulated transfer efficiency vs. beam separation, at different total energy of the particle.

Its general features agree well with the qualitative explanation given earlier. At small beam separation, significant transfer only happens at high particle energies. At beam separations around w_0 , the transfer probability is more equally spread over all possible energies, and generally much higher than in the case of small beam separation. Beam separations exceeding w_0 show no transfer at low particle energies, due to the potential barrier between the beams; the excluded energy range grows with increasing separation.

Finally, Fig. 5 shows the simulated transfer probability vs. beam separation for a thermal distribution of atoms; this diagram must be compared with the experimental result of Fig. 1. It was calculated from the data of Fig. 4 by weighting them with a thermal distribution of particle energy in a 3-dimensional harmonic oscillator. The temperature was taken to be 10% of the trap depth, which we determined in independent measurements, and which has also been consistently found in several other experiments [14].

The simulated transfer probability is maximal around a beam separation of w_0 , with a pronounced minimum at the position of perfect crossing of the beams. For separations larger than w_0 , the probability falls off steeply. Therefore, the main features of the measurement are already well reflected in these simple simulations, using a static trap potential of Gaussian beams. The largest deviation between experiment and simulation is observed for small beam separation, where the measured transfer probability is significantly higher. We attribute this

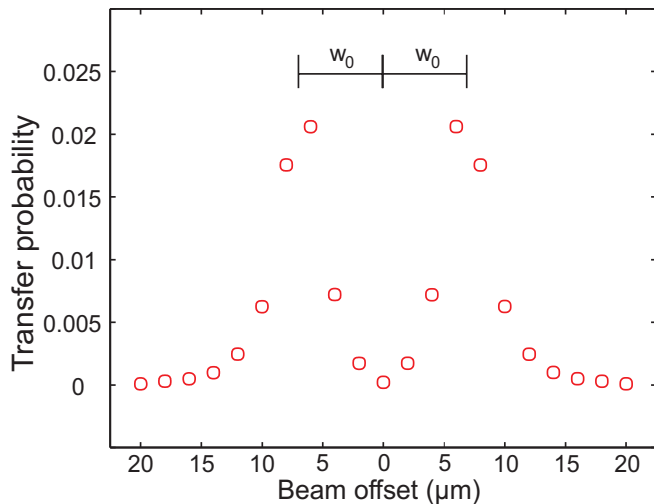


FIG. 5: Simulated transfer efficiency vs. beam separation, for a thermal distribution of particle energies as in the experiment. This diagram must be compared to Fig. 1.

to the transient situation of non-equal intensities of the beams during ramp-up and ramp-down in the experiment, which is less symmetric and may therefore mix the motion more efficiently. Other effects, such as elastic collisions between atoms in the crossing region, may also play a role in redistributing energy among the three modes of oscillation, and thus enhance the transfer efficiency at small beam separation for a cloud of atom, as in

the experiment, relative to the simulated single-particle case. We exclude, on the other hand, that experimental inaccuracies, such as non-Gaussian beam shapes or beam pointing instabilities, have broadened the central part of the measured efficiency curve, because no such broadening is observed on the outer wings of the curve.

V. CONCLUSIONS

In summary, we have measured the transfer of trapped atoms between two crossed laser beams, varying the distance between the beams. We find optimum transfer at non-zero beam separation, when the mixing between the motional degrees of freedom is favoured by the anharmonic potential, while no potential barrier is formed between the beams. The main characteristics of the experimental observations are reproduced in numerical simulations of the dynamics. They become visible already in the situation of two crossed beams of constant, equal power. Our results are relevant for designing efficient loading mechanisms for optical traps and optical lattices, and they may find applications in optical tweezer technology for nano- and micro-particles.

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